



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:11 PM UTC

PDB ID : 9N6A / pdb_00009n6a
Title : GP23 C-terminal receptor binding domain from TM4 phage
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Deposited on : 2025-02-05
Resolution : 2.91 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	NOT EXECUTED
Xtriage (Phenix)	:	2.0
EDS	:	NOT EXECUTED
Buster-report	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

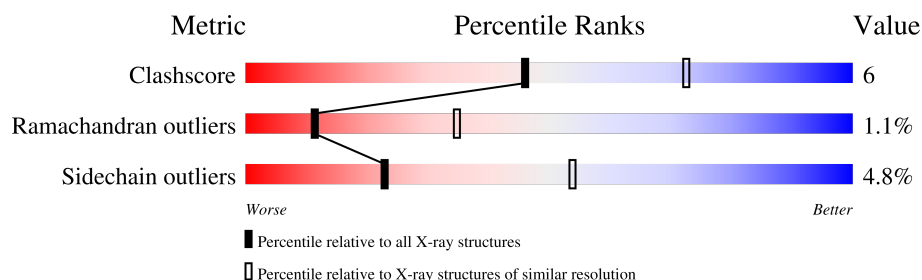
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	3213 (2.94-2.90)
Ramachandran outliers	187476	3128 (2.94-2.90)
Sidechain outliers	187428	3130 (2.94-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	627	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2765 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Green fluorescent protein, Minor tail protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2714	1718	468	520	8			

There are 31 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-12	MET	-	initiating methionine	UNP A0A059PIQ0
A	-11	LYS	-	expression tag	UNP A0A059PIQ0
A	-10	SER	-	expression tag	UNP A0A059PIQ0
A	-9	SER	-	expression tag	UNP A0A059PIQ0
A	-8	HIS	-	expression tag	UNP A0A059PIQ0
A	-7	HIS	-	expression tag	UNP A0A059PIQ0
A	-6	HIS	-	expression tag	UNP A0A059PIQ0
A	-5	HIS	-	expression tag	UNP A0A059PIQ0
A	-4	HIS	-	expression tag	UNP A0A059PIQ0
A	-3	HIS	-	expression tag	UNP A0A059PIQ0
A	-2	GLY	-	expression tag	UNP A0A059PIQ0
A	-1	SER	-	expression tag	UNP A0A059PIQ0
A	0	SER	-	expression tag	UNP A0A059PIQ0
A	1	VAL	MET	conflict	UNP A0A059PIQ0
A	2	SER	ARG	conflict	UNP A0A059PIQ0
A	30	ARG	SER	conflict	UNP A0A059PIQ0
A	72	SER	ALA	conflict	UNP A0A059PIQ0
A	206	LYS	ALA	conflict	UNP A0A059PIQ0
A	231	LEU	HIS	conflict	UNP A0A059PIQ0
A	239	GLY	-	linker	UNP A0A059PIQ0
A	240	ILE	-	linker	UNP A0A059PIQ0
A	241	GLU	-	linker	UNP A0A059PIQ0
A	242	GLU	-	linker	UNP A0A059PIQ0
A	243	ASN	-	linker	UNP A0A059PIQ0
A	244	LEU	-	linker	UNP A0A059PIQ0
A	245	TYR	-	linker	UNP A0A059PIQ0
A	246	PHE	-	linker	UNP A0A059PIQ0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	247	GLN	-	linker	UNP A0A059PIQ0
A	248	SER	-	linker	UNP A0A059PIQ0
A	249	ASN	-	linker	UNP A0A059PIQ0
A	250	ALA	-	linker	UNP A0A059PIQ0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



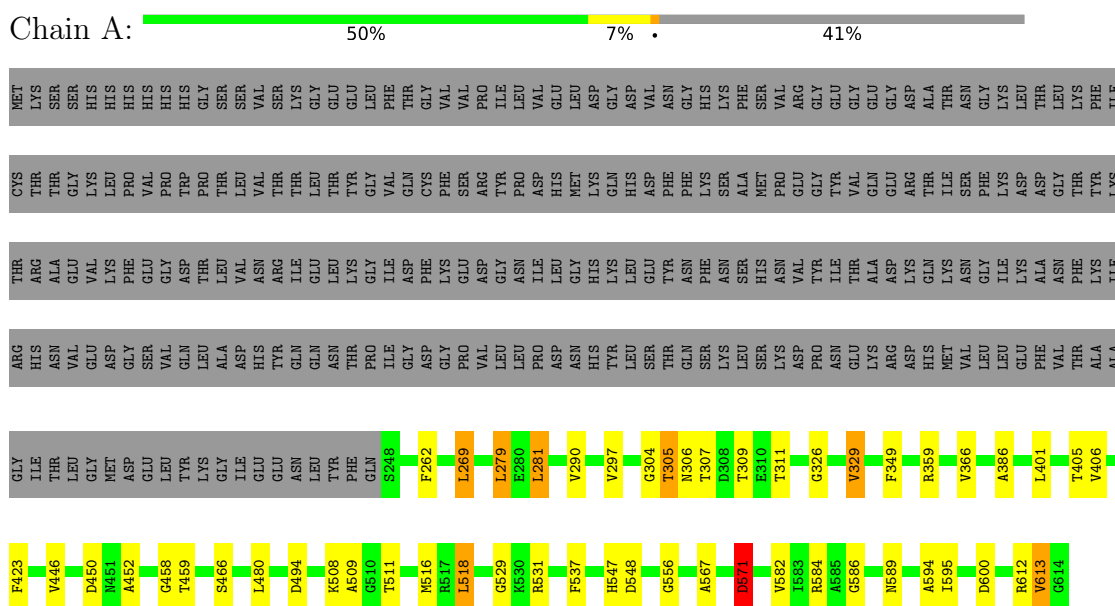
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		

Note EDS was not executed.

- Molecule 1: Green fluorescent protein, Minor tail protein



4 Data and refinement statistics

EDS was not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	125.73Å 125.73Å 97.51Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.32 – 2.91	Depositor
% Data completeness (in resolution range)	98.3 (36.32-2.91)	Depositor
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.230 , 0.280	Depositor
Wilson B-factor (Å ²)	86.2	Xtriage
Anisotropy	0.199	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.042 for h,-h-k,-l	Xtriage
Total number of atoms	2765	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.46	0/2783	0.73	0/3794

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	359	ARG	Sidechain

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2714	0	2596	31	0
2	A	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	41	0	0	8	0
All	All	2765	0	2596	31	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (31) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:LEU:HD11	3:A:814:HOH:O	1.52	1.07
1:A:589:ASN:HD22	1:A:595:ILE:HD11	1.49	0.78
1:A:458:GLY:N	3:A:802:HOH:O	2.29	0.64
1:A:571:ASP:HB3	3:A:822:HOH:O	1.98	0.63
1:A:329:VAL:HG12	1:A:349:PHE:HB3	1.81	0.61
1:A:304:GLY:O	1:A:306:ASN:N	2.33	0.60
1:A:547:HIS:HB3	3:A:813:HOH:O	2.04	0.58
1:A:279:LEU:HD11	1:A:297:VAL:HG11	1.87	0.57
1:A:594:ALA:C	1:A:595:ILE:HD13	2.31	0.56
1:A:311:THR:HG21	1:A:452:ALA:HB2	1.87	0.56
1:A:423:PHE:HB2	3:A:801:HOH:O	2.07	0.54
1:A:459:THR:HG23	1:A:494:ASP:OD1	2.07	0.54
1:A:262:PHE:HA	1:A:281:LEU:HD11	1.89	0.54
1:A:304:GLY:N	1:A:307:THR:OG1	2.41	0.53
1:A:567:ALA:N	3:A:804:HOH:O	2.44	0.51
1:A:386:ALA:HB3	1:A:401:LEU:HB2	1.94	0.50
1:A:311:THR:HG22	1:A:450:ASP:OD1	2.13	0.48
1:A:401:LEU:HD21	1:A:406:VAL:HG22	1.96	0.47
1:A:508:LYS:HG2	1:A:612:ARG:NH1	2.30	0.46
1:A:516:MET:HE2	1:A:518:LEU:HD11	1.97	0.45
1:A:529:GLY:HA2	3:A:809:HOH:O	2.16	0.45
1:A:466:SER:HA	1:A:600:ASP:O	2.16	0.45
1:A:279:LEU:CD1	1:A:297:VAL:HG11	2.48	0.43
1:A:480:LEU:HD22	1:A:586:GLY:HA2	2.01	0.43
1:A:531:ARG:NH2	1:A:548:ASP:OD2	2.50	0.43
1:A:589:ASN:ND2	1:A:595:ILE:HD11	2.24	0.42
1:A:509:ALA:O	1:A:612:ARG:NH1	2.48	0.42
1:A:309:THR:HB	1:A:450:ASP:OD1	2.19	0.41
1:A:537:PHE:HB2	1:A:584:ARG:HB3	2.02	0.41
1:A:423:PHE:O	3:A:801:HOH:O	2.21	0.41
1:A:511:THR:HB	1:A:613:VAL:HG22	2.02	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	365/627 (58%)	340 (93%)	21 (6%)	4 (1%)	11 35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	305	THR
1	A	571	ASP
1	A	326	GLY
1	A	556	GLY

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	270/498 (54%)	257 (95%)	13 (5%)	23 54

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	269	LEU
1	A	279	LEU
1	A	281	LEU
1	A	290	VAL

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Mol	Chain	Res	Type
1	A	305	THR
1	A	329	VAL
1	A	366	VAL
1	A	405	THR
1	A	446	VAL
1	A	518	LEU
1	A	571	ASP
1	A	582	VAL
1	A	613	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

Mogul was not executed - this section is therefore empty.

5.5 Carbohydrates [i](#)

Mogul was not executed - this section is therefore empty.

5.6 Ligand geometry [i](#)

Mogul was not executed - this section is therefore empty.

5.7 Other polymers [i](#)

Mogul was not executed - this section is therefore empty.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.